

IN THE CLAIMS:

Claims 1-4 (**cancelled**).

Claim 5 (**previously presented**): A compound as claimed in claim 18 wherein Zb is -O-.

Claim 6 (**previously presented**): A compound as claimed in claim 18 wherein R^{2a} is methoxy.

Claim 7 (**previously presented**): A compound as claimed in claim 18 wherein ring C is a 5-membered heteroaromatic moiety which contains 1-3 heteroatoms selected independently from O, N and S.

Claim 8 (**previously presented**): A compound as claimed in claim 18 wherein R¹ is a phenyl group or a 5-6-membered heteroaromatic group with 1-3 heteroatoms, selected independently from O, S and N, (linked via a ring carbon atom), which phenyl or heteroaromatic group is optionally substituted as defined in claim 18.

Claim 9 (**currently amended**): A compound as claimed in claim 18 wherein R² represents hydroxy, ~~halogeno~~, nitro, trifluoromethyl, C₁₋₃alkyl, cyano, amino or R⁵X¹- wherein X¹ is as defined in claim 18 and R⁵ is selected from one of the following eighteen groups:

- 1) C₁₋₄alkyl which may be unsubstituted or substituted with one or more fluorine atoms, or C₂₋₄alkyl which may be unsubstituted or substituted with 1 or 2 groups selected from hydroxy and amino;
- 2) C₂₋₃alkylX²COR¹¹ (wherein X² is as defined in claim 18 and R¹¹ represents -NR¹³R¹⁴ or -OR¹⁵ (wherein R¹³, R¹⁴ and R¹⁵ which may be the same or different are each C₁₋₂alkyl or C₁₋₂alkoxyethyl));

- 3) $C_{2-4}alkylX^3R^{16}$ (wherein X^3 is as defined in claim 18 and R^{16} is a group selected from $C_{1-3}alkyl$, cyclopentyl, cyclohexyl, pyrrolidinyl and piperidinyl which group is linked to X^3 through a carbon atom and which $C_{1-3}alkyl$ group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno and $C_{1-2}alkoxy$ and which cyclopentyl, cyclohexyl, pyrrolidinyl or piperidinyl group may carry one substituent selected from oxo, hydroxy, halogeno, $C_{1-2}alkyl$, $C_{1-2}hydroxyalkyl$ and $C_{1-2}alkoxy$);
- 4) $C_{2-3}alkylX^4C_{2-3}alkylX^5R^{22}$ (wherein X^4 and X^5 are as defined in claim 18 and R^{22} represents hydrogen or $C_{1-3}alkyl$);
- 5) $C_{1-4}alkylR^{59}$ (wherein R^{59} is a group selected from pyrrolidinyl, piperazinyl, piperidinyl, 1,3-dioxolan-2-yl, 1,3-dioxan-2-yl, 1,3-dithiolan-2-yl and 1,3-dithian-2-yl, which group is linked to $C_{1-4}alkyl$ through a carbon atom and which group may carry 1 or 2 substituents selected from oxo, hydroxy, halogeno, $C_{1-3}alkyl$, $C_{1-3}hydroxyalkyl$, $C_{1-3}alkoxy$, $C_{1-2}alkoxyC_{1-3}alkyl$ and $C_{1-2}alkylsulphonylC_{1-3}alkyl$) or $C_{2-4}alkylR^{60}$ (wherein R^{60} is a group selected from morpholino, thiomorpholino, pyrrolidin-1-yl, piperazin-1-yl and piperidino which group may carry 1 or 2 substituents selected from oxo, hydroxy, halogeno, $C_{1-3}alkyl$, $C_{1-3}hydroxyalkyl$, $C_{1-3}alkoxy$, $C_{1-2}alkoxyC_{1-3}alkyl$ and $C_{1-2}alkylsulphonylC_{1-3}alkyl$);
- 6) $C_{3-4}alkenylR^{61}$ (wherein R^{61} represents R^{59} or R^{60} as defined herein);
- 7) $C_{3-4}alkynylR^{61}$ (wherein R^{61} represents R^{59} or R^{60} as defined herein);
- 8) R^{29} (wherein R^{29} is as defined in claim 18);
- 9) $C_{1-4}alkylR^{29}$ (wherein R^{29} is as defined in claim 18);
- 10) 1- R^{29} prop-1-en-3-yl or 1- R^{29} but-2-en-4-yl (wherein R^{29} is as defined in claim 18 with the proviso that when R^5 is 1- R^{29} prop-1-en-3-yl, R^{29} is linked to the alkenyl group via a carbon atom);
- 11) 1- R^{29} prop-1-yn-3-yl or 1- R^{29} but-2-yn-4-yl (wherein R^{29} is as defined in claim 18 with the proviso that when R^5 is 1- R^{29} prop-1-yn-3-yl, R^{29} is linked to the alkynyl group via a carbon atom);
- 12) $C_{1-5}alkylX^6R^{29}$ (wherein X^6 and R^{29} are as defined in claim 18);
- 13) 1-($R^{29}X^7$)but-2-en-4-yl (wherein X^7 and R^{29} are as defined in claim 18);

- 14) 1-(R²⁹X⁸)but-2-yn-4-yl (wherein X⁸ and R²⁹ are as defined in claim 18);
 15) C₂₋₃alkylX⁹C₁₋₂alkylR²⁹ (wherein X⁹ and R²⁹ are as defined in claim 18);
 16) R²⁸ (wherein R²⁸ is as defined in claim 18);
 17) C₂₋₃alkylX⁹C₁₋₂alkylR²⁸ (wherein X⁹ and R²⁸ are as defined in claim 18); and
 18) C₂₋₃alkylR⁵⁴C₁₋₂alkylX⁹R⁵⁵ (wherein X⁹, R⁵⁴ and R⁵⁵ are as defined in claim 18);
 and additionally wherein any C₁₋₅alkyl, C₂₋₅alkenyl or C₂₋₅alkynyl group in R⁵X¹- may bear one or more substituents selected from hydroxy, halogeno and amino.

Claim 10 (**previously presented**): A compound as claimed in claim 18 wherein R² represents 2-methoxyethoxy, 2-(2-methoxyethoxy)ethoxy, 3-methoxypropoxy, 2-methylsulphonylethoxy, 3-methylsulphonylpropoxy, 2-(tetrahydropyran-4-yloxy)ethoxy, 3-(tetrahydropyran-4-yloxy)propoxy, 2-(4-methylpiperazin-1-yl)ethoxy, 3-(4-methylpiperazin-1-yl)propoxy, 2-morpholinoethoxy, 3-morpholinopropoxy, 2-(imidazol-1-yl)ethoxy, 3-(imidazol-1-yl)propoxy 2-(1,1-dioxothiomorpholino)ethoxy, 3-(1,1-dioxothiomorpholino)propoxy, 2-(1,2,3-triazol-1-yl)ethoxy, 3-(1,2,3-triazol-1-yl)propoxy, 2-(N-methoxyacetyl-N-methylamino)ethoxy, 3-(N-methoxyacetyl-N-methylamino)propoxy, N-methylpiperidin-3-ylmethoxy, 4-(pyrrolidin-1-yl)but-2-en-yloxy, 2-(2-oxopyrrolidin-1-yl)ethoxy, 3-(2-oxopyrrolidin-1-yl)propoxy, 2-(pyrrolidin-1-yl)ethoxy, 3-(pyrrolidin-1-yl)propoxy, 2-(2-(pyrrolidin-1-yl)ethoxy)ethoxy, 2-(2-(4-methylpiperazin-1-yl)ethoxy)ethoxy, 2-piperidinoethoxy, 3-piperidinopropoxy, 2-(methylpiperidino)ethoxy, 3-(methylpiperidino)propoxy, 2-(ethylpiperidino)ethoxy, 3-(ethylpiperidino)propoxy, 2-((2-methoxyethyl)piperidino)ethoxy, 3-((2-methoxyethyl)piperidino)propoxy, 2-((2-methylsulphonyl)ethylpiperidino)ethoxy, 3-((2-methylsulphonyl)ethylpiperidino)propoxy, piperidin-3-ylmethoxy, piperidin-4-ylmethoxy, 2-(piperidin-3-yl)ethoxy, 2-(piperidin-4-yl)ethoxy, 3-(piperidin-3-yl)propoxy, 3-(piperidin-4-yl)propoxy, 2-(methylpiperidin-3-yl)ethoxy, 2-(methylpiperidin-4-yl)ethoxy, 3-(methylpiperidin-3-yl)propoxy, 3-(methylpiperidin-4-yl)propoxy, 2-(ethylpiperidin-3-yl)ethoxy,

2-(ethylpiperidin-4-yl)ethoxy, 3-(ethylpiperidin-3-yl)propoxy,
3-(ethylpiperidin-4-yl)propoxy, 2-((2-methoxyethyl)piperidin-3-yl)ethoxy,
2-((2-methoxyethyl)piperidin-4-yl)ethoxy, 3-((2-methoxyethyl)piperidin-3-yl)propoxy,
3-((2-methoxyethyl)piperidin-4-yl)propoxy,
2-((2-methylsulphonyl)ethyl)piperidin-3-yl)ethoxy,
2-((2-methylsulphonyl)ethyl)piperidin-4-yl)ethoxy,
3-((2-methylsulphonyl)ethyl)piperidin-3-yl)propoxy,
3-((2-methylsulphonyl)ethyl)piperidin-4-yl)propoxy, 1-isopropylpiperidin-2-ylmethyl,
1-isopropylpiperidin-3-ylmethyl, 1-isopropylpiperidin-4-ylmethyl,
2-(1-isopropylpiperidin-2-yl)ethyl, 2-(1-isopropylpiperidin-3-yl)ethyl,
2-(1-isopropylpiperidin-4-yl)ethyl, 3-(1-isopropylpiperidin-2-yl)propyl,
3-(1-isopropylpiperidin-3-yl)propyl, 3-(1-isopropylpiperidin-4-yl)propyl,
3-(4-methylpiperazin-1-yl)propoxy, 1-methylpiperidin-4-ylmethoxy,
1-(2-methylsulphonyl)ethyl)piperidin-4-ylmethoxy,
1-(2-pyrrolidinylethyl)piperidin-4-ylmethoxy,
1-(3-pyrrolidinylpropyl)piperidin-4-ylmethoxy, 1-(2-piperidinylethyl)piperidin-4-ylmethoxy,
1-(3-piperidinylpropyl)piperidin-4-ylmethoxy, 1-(2-morpholinoethyl)piperidin-4-ylmethoxy,
1-(3-morpholinopropyl)piperidin-4-ylmethoxy,
1-(2-thiomorpholinoethyl)piperidin-4-ylmethoxy,
1-(3-thiomorpholinopropyl)piperidin-4-ylmethoxy,
1-(2-azetidinyethyl)piperidin-4-ylmethoxy or 1-(3-azetidinypropyl)piperidin-4-ylmethoxy.

Claim 11 (**previously presented**): A compound as claimed in claim 18 selected from:
4-(5-(4-methoxyphenyl)pyrazol-3-yloxy)-6-methoxy-7-(1-methylpiperidin-4-ylmethoxy)-
quinazoline,
4-(5-(4-methoxyphenyl)pyrazol-3-yloxy)-6-methoxy-7-(3-(4-methylpiperazin-1-yl)-
propoxy)quinazoline,
6-methoxy-7-(2-(2-methoxyethoxy)ethoxy)-4-(5-phenylpyrazol-3-yloxy)quinazoline,
4-(5-(3-furyl)pyrazol-3-yloxy)-6-methoxy-7-(3-morpholinopropoxy)quinazoline,

6-methoxy-7-(3-morpholinopropoxy)-4-(5-phenylpyrazol-3-yloxy)quinazoline,
7-(2-(imidazol-1-yl)ethoxy)-6-methoxy-4-(5-phenylpyrazol-3-yloxy)quinazoline,
4-(5-(4-chlorophenyl)pyrazol-3-yloxy)-6-methoxy-7-(3-morpholinopropoxy)quinazoline,
6-methoxy-7-(3-(4-methylpiperazin-1-yl)propoxy)-4-(5-phenylpyrazol-3-yloxy)-quinazoline,
6-methoxy-7-(2-methoxyethoxy)-4-(5-phenylpyrazol-3-yloxy)quinazoline,
4-(5-(4-methoxyphenyl)pyrazol-3-yloxy)-6-methoxy-7-(2-(1,2,3-triazol-1-yl)ethoxy)-
quinazoline and
4-(5-(4-methoxyphenyl)pyrazol-3-yloxy)-6-methoxy-7-(1-(2-methylsulphonylethyl)-
piperidin-4-ylmethoxy)quinazoline,
and salts thereof.

Claim 12 (**previously presented**): A compound as claimed in claim 18 selected from:

7-(2-methoxyethoxy)-4-(5-phenylpyrazol-3-yloxy)quinazoline,
4-(5-(2-fluorophenyl)pyrazol-3-yloxy)-6-methoxy-7-(3-morpholinopropoxy)quinazoline,
6-methoxy-7-(3-morpholinopropoxy)-4-(5-(3-nitrophenyl)pyrazol-3-yloxy)quinazoline,
6-methoxy-7-(3-morpholinopropoxy)-4-(5-(4-nitrophenyl)pyrazol-3-yloxy)quinazoline,
6-methoxy-7-(3-morpholinopropoxy)-4-(5-(4-pyridyl)pyrazol-3-yloxy)quinazoline,
4-(5-(4-fluorophenyl)pyrazol-3-yloxy)-6-methoxy-7-(3-morpholinopropoxy)quinazoline, and
6-methoxy-7-(2-methoxyethoxy)-4-(5-(4-methoxyphenyl)pyrazol-3-yloxy)quinazoline,
and salts thereof.

Claim 13 (**previously presented**): A method for producing an antiangiogenic and/or vascular permeability reducing effect in a warm-blooded animal in need of such treatment which comprises administering to such animal an effective amount of a compound selected from the group consisting of:

6-methoxy-7-(1-methylpiperidin-4-ylmethoxy)-4-(5-phenylpyrazol-3-ylamino)-quinazoline
and
6,7-dimethoxy-4-(5-phenylpyrazol-3-yloxy)quinazoline
and pharmaceutically acceptable salts thereof.

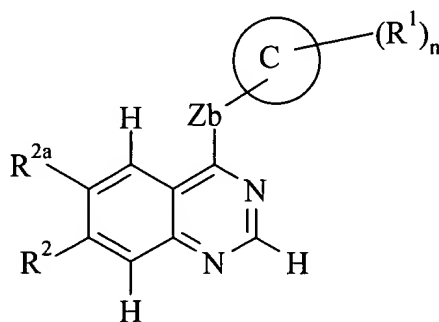
Claim 14 (**previously presented**): A compound as claimed in any one of claims 18 and 5 to 12 in the form of a pharmaceutically acceptable salt.

Claim 15 (**cancelled**).

Claim 16 (**previously presented**): A pharmaceutical composition which comprises as active ingredient a compound of formula II or a pharmaceutically acceptable salt thereof as claimed in any one of claims 18 and 5 to 12 in association with a pharmaceutically acceptable excipient or carrier.

Claim 17 (**previously presented**): A method for producing an antiangiogenic and/or vascular permeability reducing effect in a warm-blooded animal in need of such treatment which comprises administering to said animal an effective amount of a compound of formula II as defined in any one of claims 18 and 5 to 12 or a pharmaceutically acceptable salt thereof.

Claim 18 (**currently amended**): A compound of the formula II:



II

wherein:

ring C is a 5-6-membered heterocyclic moiety which may be saturated or unsaturated, which may be aromatic or non-aromatic, and which contains 1-3 heteroatoms selected independently from O, N and S;

Zb is -O- or -S-;

R¹ represents hydrogen, C₁₋₄alkyl, C₁₋₄alkoxymethyl, di(C₁₋₄alkoxy)methyl, C₁₋₄alkanoyl, trifluoromethyl, cyano, amino, C₂₋₅alkenyl, C₂₋₅alkynyl, a phenyl group, a benzyl group or a 5-6-membered heterocyclic group with 1-3 heteroatoms, selected independently from O, S and N, which heterocyclic group may be aromatic or non-aromatic and may be saturated (~~linked and linked~~ via a ring carbon or nitrogen ~~atom, atom~~) or unsaturated (~~linked and linked~~ via a ring carbon ~~atom, atom~~), and which phenyl, benzyl or heterocyclic group may bear on one or more ring carbon atoms up to 5 substituents selected from hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro, C₂₋₄alkanoyl, C₁₋₄alkanoylamino, C₁₋₄alkoxycarbonyl, C₁₋₄alkylsulphanyl, C₁₋₄alkylsulphinyl, C₁₋₄alkylsulphonyl, carbamoyl, N-C₁₋₄alkylcarbamoyl, N,N-di(C₁₋₄alkyl)carbamoyl, aminosulphonyl, N-C₁₋₄alkylaminosulphonyl, N,N-di(C₁₋₄alkyl)aminosulphonyl, C₁₋₄alkylsulphonylamino, C₁₋₄alkylamino, C₁₋₄hydroxyalkyl, C₁₋₄aminoalkyl, C₁₋₄haloalkyl, C₁₋₄hydroxyalkoxy, carboxy and a saturated heterocyclic group selected from morpholino, thiomorpholino, pyrrolidinyl, piperazinyl, piperidinyl imidazolidinyl and pyrazolidinyl, which saturated heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₃alkyl, C₁₋₃alkoxy, C₁₋₃alkanoyloxy, trifluoromethyl, cyano, amino, nitro and C₁₋₄alkoxycarbonyl; and additionally R¹ may represent carboxy, C₃₋₇cycloalkyl, C₃₋₇cycloalkylC₁₋₃alkyl, or phenylC₂₋₄alkyl wherein the phenyl moiety may bear up to 5 substituents selected from the list herein defined for a phenyl ring which is directly linked to ring C;

n is an integer from 0 to 5;

m is an integer from 0 to 3;

R^2 represents hydroxy, cyano, nitro, trifluoromethyl, C_{1-3} alkylsulphanyl, $-NR^3R^4$ (~~wherein, wherein~~ R^3 and R^4 , which may be the same or different, each represents hydrogen or C_{1-3} alkyl), C_{1-3} alkyl,

or R^2 represents R^5X^1 (~~wherein, wherein~~ X^1 represents a direct bond, -O-, $-CH_2-$, $-OCO-$, carbonyl, -S-, -SO-, $-SO_2-$, $-NR^6CO-$, $-CONR^7-$, $-SO_2NR^8-$, $-NR^9SO_2-$ or $-NR^{10}-$ (~~wherein, wherein~~ R^6 , R^7 , R^8 , R^9 and R^{10} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl), ~~and, and~~ R^5 is selected from one of the following eighteen groups:

- 1) hydrogen or C_{1-5} alkyl which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, fluoro, chloro, bromo and amino;
- 2) $C_{1-5}alkylX^2COR^{11}$ (~~wherein, wherein~~ X^2 represents -O- or $-NR^{12}-$ (~~in which, in which~~ R^{12} represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl) ~~and, and~~ R^{11} represents C_{1-3} alkyl, $-NR^{13}R^{14}$ or $-OR^{15}$ (~~wherein, wherein~~ R^{13} , R^{14} and R^{15} which may be the same or different each represents hydrogen, C_{1-3} alkyl, C_{4-5} alkyl or C_{1-3} alkoxy C_{2-3} alkyl));
- 3) $C_{1-5}alkylX^3R^{16}$ (~~wherein, wherein~~ X^3 represents -O-, -S-, -SO-, $-SO_2-$, $-OCO-$, $-NR^{17}CO-$, $-CONR^{18}-$, $-SO_2NR^{19}-$, $-NR^{20}SO_2-$ or $-NR^{21}-$ (~~wherein, wherein~~ R^{17} , R^{18} , R^{19} , R^{20} and R^{21} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl) ~~and, and~~ R^{16} represents hydrogen, C_{1-3} alkyl, cyclopentyl, cyclohexyl or a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which C_{1-3} alkyl group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno and C_{1-4} alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-4} alkyl, C_{1-4} hydroxyalkyl, C_{1-4} alkoxy, C_{1-4} cyanoalkyl and C_{1-4} alkoxycarbonyl) ~~C_{1-4} alkoxycarbonyl~~);
- 4) $C_{1-5}alkylX^4C_{1-5}alkylX^5R^{22}$ (~~wherein, wherein~~ X^4 and X^5 ~~which, which~~ may be the same or different ~~are, are~~ each -O-, -S-, -SO-, $-SO_2-$, $-NR^{23}CO-$, $-CONR^{24}-$, $-SO_2NR^{25}-$, $-NR^{26}SO_2-$ or $-NR^{27}-$ (~~wherein, wherein~~ R^{23} , R^{24} , R^{25} , R^{26} and R^{27} each

- independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) ~~and, and~~ and R²² represents hydrogen, C₁₋₃alkyl or ~~C₁₋₃alkoxyC₂₋₃alkyl~~ C₁₋₃alkoxyC₂₋₃alkyl;
- 5) R²⁸ ~~(wherein, wherein R²⁸ is a 5-6-membered saturated heterocyclic group (linked, linked via carbon or nitrogen) with, with 1-2 heteroatoms, selected selected~~ independently from O, S and N, which heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₄alkyl, C₁₋₄hydroxyalkyl, C₁₋₄alkoxy, C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl and ~~C₁₋₄alkoxycarbonyl~~) C₁₋₄alkoxycarbonyl;
- 6) C₁₋₅alkylR²⁸ ~~(wherein, wherein R²⁸ is as defined herein) herein;~~ herein;
- 7) C₂₋₅alkenylR²⁸ ~~(wherein, wherein R²⁸ is as defined herein) herein;~~ herein;
- 8) C₂₋₅alkynylR²⁸ ~~(wherein, wherein R²⁸ is as defined herein) herein;~~ herein;
- 9) R²⁹ ~~(wherein, wherein R²⁹ represents a pyridone group, a phenyl group or a 5-6-membered aromatic heterocyclic group (linked, linked via carbon or nitrogen) with, with 1-3 heteroatoms selected from O, N and S, which pyridone, phenyl or aromatic heterocyclic group may carry up to 5 substituents on an available carbon atom selected from hydroxy, halogeno, amino, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄hydroxyalkyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, C₁₋₄hydroxyalkoxy, carboxy, trifluoromethyl, cyano, -CONR³⁰R³¹ and -NR³²COR³³ (wherein, wherein R³⁰, R³¹, R³² and R³³, which may be the same or different, each represents hydrogen, C₁₋₄alkyl or C₁₋₃alkoxyC₂₋₃alkyl));~~ herein;
- 10) C₁₋₅alkylR²⁹ ~~(wherein, wherein R²⁹ is as defined herein) herein;~~ herein;
- 11) C₂₋₅alkenylR²⁹ ~~(wherein, wherein R²⁹ is as defined herein) herein;~~ herein;
- 12) C₂₋₅alkynylR²⁹ ~~(wherein, wherein R²⁹ is as defined herein) herein;~~ herein;
- 13) C₁₋₅alkylX⁶R²⁹ ~~(wherein, wherein X⁶ represents -O-, -S-, -SO-, -SO₂-, -NR³⁴CO-, -CONR³⁵-, -SO₂NR³⁶-, -NR³⁷SO₂- or -NR³⁸- (wherein, wherein R³⁴, R³⁵, R³⁶, R³⁷ and R³⁸ each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and, and R²⁹ is as defined herein) herein;~~ herein;
- 14) C₂₋₅alkenylX⁷R²⁹ ~~(wherein, wherein X⁷ represents -O-, -S-, -SO-, -SO₂-, -NR³⁹CO-, -CONR⁴⁰-, -SO₂NR⁴¹-, -NR⁴²SO₂- or -NR⁴³- (wherein, wherein R³⁹, R⁴⁰, R⁴¹, R⁴² and~~ herein;

R^{43} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl) ~~and~~,
and R^{29} is as defined ~~herein~~) herein;

15) C_{2-5} alkynyl X^8R^{29} ~~(wherein, wherein X^8 represents -O-, -S-, -SO-, -SO₂-, -NR⁴⁴CO-,
-CONR⁴⁵-, -SO₂NR⁴⁶-, -NR⁴⁷SO₂- or -NR⁴⁸- (wherein, wherein R^{44} , R^{45} , R^{46} , R^{47} and
 R^{48} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl) ~~and~~,
and R^{29} is as defined ~~herein~~) herein;~~

16) C_{1-3} alkyl X^9C_{1-3} alkyl R^{29} ~~(wherein, wherein X^9 represents -O-, -S-, -SO-, -SO₂-,
-NR⁴⁹CO-, -CONR⁵⁰-, -SO₂NR⁵¹-, -NR⁵²SO₂- or -NR⁵³- (wherein, wherein R^{49} , R^{50} ,
 R^{51} , R^{52} and R^{53} each independently represents hydrogen, C_{1-3} alkyl or
 C_{1-3} alkoxy C_{2-3} alkyl) ~~and~~, and R^{29} is as defined ~~herein~~) herein;~~

17) C_{1-3} alkyl X^9C_{1-3} alkyl R^{28} ~~(wherein, wherein X^9 and R^{28} are as defined ~~herein~~) herein;~~
and

18) C_{1-3} alkyl $R^{54}C_{1-3}$ alkyl X^9R^{55} ~~(wherein, wherein X^9 is as defined herein and R^{54} and R^{55}
are each independently selected from hydrogen, C_{1-3} alkyl, cyclopentyl, cyclohexyl
and a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected
independently from O, S and N, which C_{1-3} alkyl group may bear 1 or 2 substituents
selected from oxo, hydroxy, halogeno and C_{1-4} alkoxy and which cyclic group may
bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-4} alkyl,
 C_{1-4} hydroxyalkyl, C_{1-4} alkoxy, C_{1-4} cyanoalkyl and C_{1-4} alkoxycarbonyl) ~~with, with the~~
proviso that R^{54} cannot be hydrogen;~~

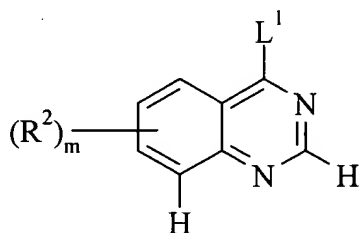
and additionally wherein any C_{1-5} alkyl, C_{2-5} alkenyl or C_{2-5} alkynyl group in R^5X^1 - may
bear one or more substituents selected from hydroxy, halogeno and amino; provided that
 R^2 is not hydrogen, substituted or unsubstituted C_{1-5} alkyl, C_{1-5} alkoxy, phenoxy or
phenyl C_{1-5} alkoxy; and

R^{2a} represents hydrogen, halogeno, C_{1-3} alkyl, C_{1-3} alkoxy, C_{1-3} alkylthio, -NR^{3a}R^{4a} ~~(wherein,
wherein R^{3a} and R^{4a} , which may be the same or different, each represents hydrogen or
 C_{1-3} alkyl) ~~or, or $R^{5a}(CH_2)_{za}X^{1a}$ (wherein, wherein R^{5a} is a 5- or 6-membered saturated
heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which
heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno,~~~~

C₁₋₄alkyl, C₁₋₄hydroxyalkyl and C₁₋₄alkoxy, z_a is an integer from 0 to 4 and X^{1a} represents a direct bond, -O-, -CH₂-, -S-, -SO-, -SO₂-, -NR^{6a}CO-, -CONR^{7a}-, -SO₂NR^{8a}-, -NR^{9a}SO₂- or -NR^{10a}- (wherein, wherein R^{6a}, R^{7a}, R^{8a}, R^{9a} and R^{10a} each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl);
or a salt thereof.

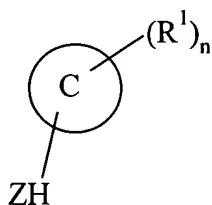
19 (previously presented): A process for the preparation of a compound of formula II or salt thereof, as defined in claim 18, which comprises:

(a) the reaction of a compound of the formula III:



(III)

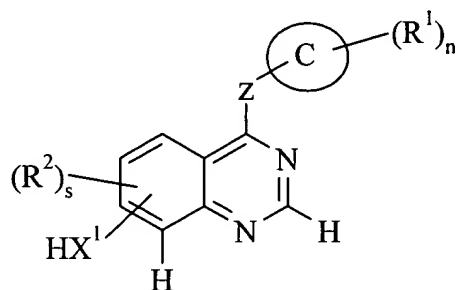
(wherein R² and m are as defined in claim 18 and L¹ is a displaceable moiety), with a compound of the formula IV:



(IV)

(wherein ring C, R¹, Z and n are as defined in claim 18);

- (b) compounds of formula II and salts thereof wherein at least one R^2 is R^5X^1 wherein R^5 is as defined in claim 18 and X^1 is -O-, -S-, -OCO- or -NR¹⁰- (wherein R^{10} independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) may be prepared by the reaction of a compound of the formula V:



(V)

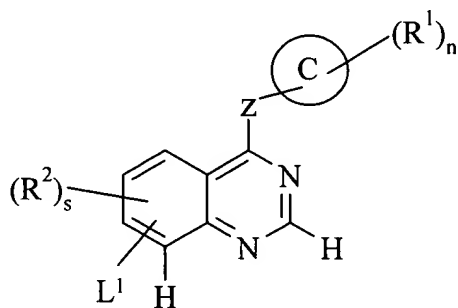
(wherein ring C, Z, R^1 , R^2 and n are as defined in claim 18 and X^1 is as defined herein in this section and s is an integer from 0 to 2) with a compound of formula VI:



(VI)

(wherein R^5 is as defined in claim 18 and L^1 is as defined herein);

- (c) compounds of the formula II and salts thereof wherein at least one R^2 is R^5X^1 wherein R^5 is as defined in claim 18 and X^1 is -O-, -S-, -OCO- or -NR¹⁰- (wherein R^{10} represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) may be prepared by the reaction of a compound of the formula VII:



(VII)

with a compound of the formula VIII:



(wherein R^1 , R^2 , R^5 , ring C, Z and n are as defined in claim 18 and s and L^1 are as defined herein and X^1 is as defined herein in this section);

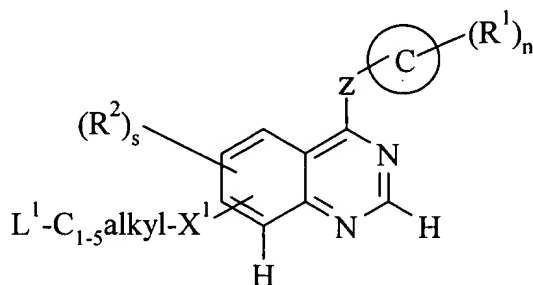
(d) compounds of the formula II and salts thereof wherein at least one R^2 is R^5X^1 wherein X^1 is as defined in claim 18 and R^5 is $C_{1-5}alkylR^{62}$, wherein R^{62} is selected from one of the following nine groups:

- 1) $X^{10}C_{1-3}alkyl$ (wherein X^{10} represents -O-, -S-, -SO₂-, -NR⁶³CO- or -NR⁶⁴SO₂- (wherein R^{63} and R^{64} which may be the same or different are each hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$);
- 2) NR⁶⁵R⁶⁶ (wherein R^{65} and R^{66} which may be the same or different are each hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$);
- 3) $X^{11}C_{1-5}alkylX^5R^{22}$ (wherein X^{11} represents -O-, -S-, -SO₂-, -NR⁶⁷CO-, -NR⁶⁸SO₂- or -NR⁶⁹- (wherein R^{67} , R^{68} , and R^{69} which may be the same or different are each hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$) and X^5 and R^{22} are as defined in claim 18);
- 4) R^{28} (wherein R^{28} is as defined in claim 18);
- 5) $X^{12}R^{29}$ (wherein X^{12} represents -O-, -S-, -SO₂-, -NR⁷⁰CO-, -NR⁷¹SO₂-, or -NR⁷²- (wherein R^{70} , R^{71} , and R^{72} which may be the same or different are each hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$) and R^{29} is as defined in claim 18);
- 6) $X^{13}C_{1-5}alkylR^{29}$, preferably $X^{13}C_{1-3}alkylR^{29}$, (wherein X^{13} represents -O-, -S-, -SO₂-, -NR⁷³CO-, -NR⁷⁴SO₂- or -NR⁷⁵- (wherein R^{73} , R^{74} and R^{75} each independently represents hydrogen, $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$) and R^{29} is as defined in claim 18);
- 7) R^{29} (wherein R^{29} is as defined in claim 18);

8) $X^{14}C_{1-3}alkylR^{28}$ (wherein X^{14} represents -O-, -S-, -SO₂-, -NR⁷⁶CO-, -NR⁷⁷SO₂- or -NR⁷⁸- (wherein R⁷⁶, R⁷⁷ and R⁷⁸ each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl) and R²⁸ is as defined in claim 18); and

9) $R^{54}C_{1-3}alkylX^9R^{55}$ (wherein R⁵⁴, R⁵⁵ and X⁹ are as defined in claim 18);

may be prepared by reacting a compound of the formula IX:



(IX)

(wherein X¹, R¹, R², ring C, Z and n are as defined in claim 18 and s and L¹ are as defined herein) with a compound of the formula X:



(X)

(wherein R⁶² is as defined herein);

- (e) compounds of the formula II and salts thereof wherein one or more of the substituents (R²)_m is represented by -NR⁷⁹R⁸⁰, where one (and the other is hydrogen) or both of R⁷⁹ and R⁸⁰ are C₁₋₃alkyl, may be prepared by the reaction of compounds of formula II wherein the substituent (R²)_m is an amino group and an alkylating agent;
- (f) compounds of the formula II and salts thereof wherein X¹ is -SO- or -SO₂- may be prepared by oxidation from the corresponding compound in which X¹ is -S- or -SO-; and when a salt of a compound of formula II is required, reaction of the compound obtained with an acid or base whereby to obtain the desired salt.